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STRUCTURE FILE UPDATES: 25 AUG 2011 HIGHEST RN 1323485-64-8
DICTIONARY FILE UPDATES: 25 AUG 2011 HIGHEST RN 1323485-64-8

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REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=>

Uploading C:\Users\rdesai\Documents\STN Express 8.4\Queries\10581177.str



chain nodes :

15 16 17 18

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14

chain bonds :

8-15 16-17 17-18

ring bonds :

1-2 1-6 1-7 1-10 2-3 3-4 4-5 5-6 7-8 8-9 9-10 9-11 10-14 11-12 12-13
13-14

exact/norm bonds :

1-2 1-6 1-7 1-10 2-3 3-4 4-5 5-6 7-8 8-9 8-15 16-17 17-18

normalized bonds :

26/08/2011

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9-10 9-11 10-14 11-12 12-13 13-14

Match level :

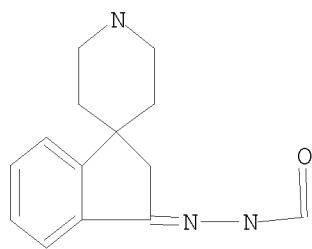
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11:Atom 12:Atom 13:Atom 14:Atom 15:CLASS 16:CLASS 17:CLASS 18:CLASS

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l1 ful

FULL SEARCH INITIATED 16:16:49 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 430 TO ITERATE

100.0% PROCESSED 430 ITERATIONS

41 ANSWERS

SEARCH TIME: 00.00.01

L2 41 SEA SSS FUL L1

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FULL SCREEN SEARCH COMPLETED - 430 TO ITERATE

100.0% PROCESSED 430 ITERATIONS

41 ANSWERS

SEARCH TIME: 00.00.01

L3 41 SEA SSS FUL L1

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This file contains CAS Registry Numbers for easy and accurate substance identification.

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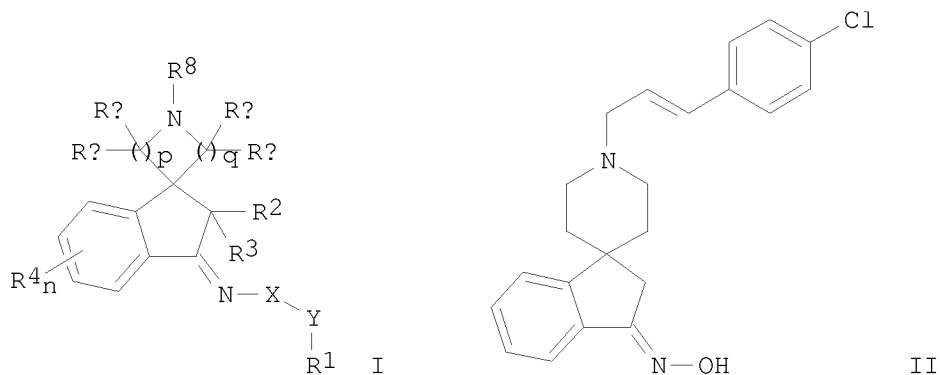
10581177.trn

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L4 1 L3

=> d abs bib hitstr

L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2011 ACS on STN
GI



AB Title compds. I [X = O, amino; Y = bond, CO, CS, SO0-2; R1 = H, alkyl, alkoxy carbonyl, etc.; R2-3 = H, halo, CN, alkyl, etc.; R4 = halo, NO2, CN, etc.; Ra = H, halo, OH, CN, etc.; p, q = 0-6; R8 = alk(en/yn)yl, etc.] are prepared For instance, II is prepared in 3 steps from spiro[indan-1-one-3,4'-piperidine]-1'-carboxylic acid tert Bu ester, 4-chlorocinnamyl chloride and hydroxylamine (E (dominant) and Z oximes isolated). Selected example compds. gave >80% control of *Spodoptera littoralis*. I are useful in controlling insects, acarines, nematodes or molluscs.

AN 2005:570877 CAPLUS

DN 143:77964

TI Preparation of insecticidal spiroindane derivatives

IN Cassayre, Jerome; Molleyres, Louis-Pierre; Maienfisch, Peter; Cederbaum, Fredrik

PA Syngenta Participations A.-G., Switz.

SO PCT Int. Appl., 114 pp.

CODEN: PIXXD2

DT Patent

LA English

FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005058836	A1	20050630	WO 2004-IB4108	20041209
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
EP 1697327	A1	20060906	EP 2004-806338	20041209
EP 1697327	B1	20110713		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,				

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IE, SI, LT, FI, RO, CY, TR, BG, CZ, EE, HU, PL, SK, IS
BR 2004017555 A 20070327 BR 2004-17555 20041209
JP 2007516253 T 20070621 JP 2006-543659 20041209
AT 516273 T 20110715 AT 2004-806338 20041209
IN 2006CN02077 A 20070706 IN 2006-CN2077 20060612
US 20080306101 A1 20081211 US 2008-581177 20080828
PRAI GB 2003-28906 A 20031212
WO 2004-IB4108 W 20041209

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

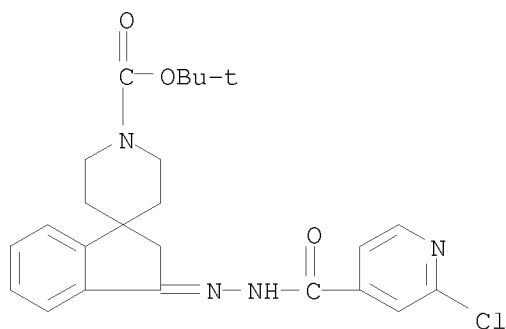
OS CASREACT 143:77964; MARPAT 143:77964

IT 855849-47-7P

RL: AGR (Agricultural use); BSU (Biological study, unclassified); RCT
(Reactant); SPN (Synthetic preparation); BIOL (Biological study); PREP
(Preparation); RACT (Reactant or reagent); USES (Uses)
(preparation of insecticidal spiroindane derivs. as insecticides,
acaracides, molluscicides and nematocides)

RN 855849-47-7 CAPLUS

CN Spiro[1H-indene-1,4'-piperidine]-1'-carboxylic acid,
3-[2-[(2-chloro-4-pyridinyl)carbonyl]hydrazinylidene]-2,3-dihydro-,
1,1-dimethylethyl ester (CA INDEX NAME)



IT 855849-48-8P 855849-49-9P 855849-50-2P
855849-51-3P 855849-52-4P 855849-53-5P
855849-54-6P 855849-55-7P 855849-56-8P
855849-57-9P 855849-58-0P 855849-59-1P
855849-60-4P 855849-61-5P 855849-62-6P
855849-63-7P 855849-64-8P 855849-65-9P
855849-66-0P 855849-67-1P 855849-68-2P
855849-69-3P 855849-70-6P 855849-71-7P
855849-72-8P 855849-73-9P 855849-74-0P
855849-75-1P 855849-76-2P 855849-77-3P
855849-78-4P 855849-79-5P 855849-80-8P
855849-81-9P 855849-85-3P 855849-86-4P
855849-87-5P 855849-94-4P

RL: AGR (Agricultural use); BSU (Biological study, unclassified); SPN
(Synthetic preparation); BIOL (Biological study); PREP (Preparation); USES
(Uses)

(preparation of insecticidal spiroindane derivs. as insecticides,
acaracides, molluscicides and nematocides)

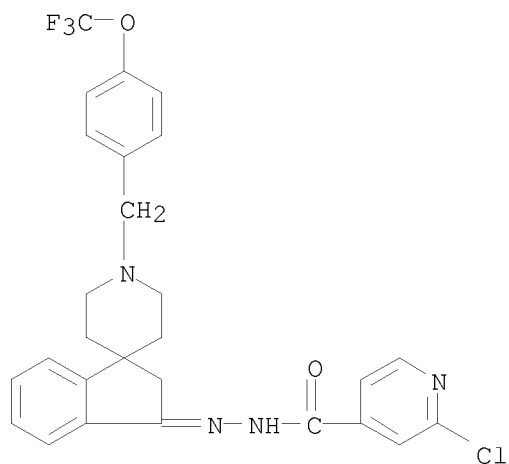
RN 855849-48-8 CAPLUS

CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[2,3-dihydro-1'-[[4-(trifluoromethoxy)phenyl]methyl]spiro[1H-indene-1,4'-

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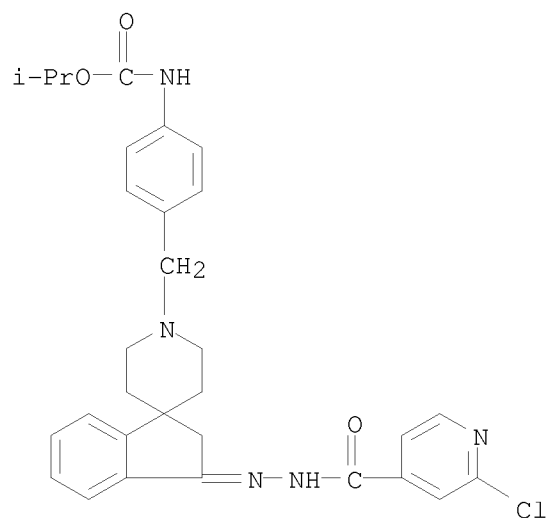
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piperidin]-3-ylidene]hydrazide (CA INDEX NAME)



RN 855849-49-9 CAPLUS

CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[2,3-dihydro-1'-[[4-[(1-methylethoxy)carbonyl]amino]phenyl)methyl]spiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

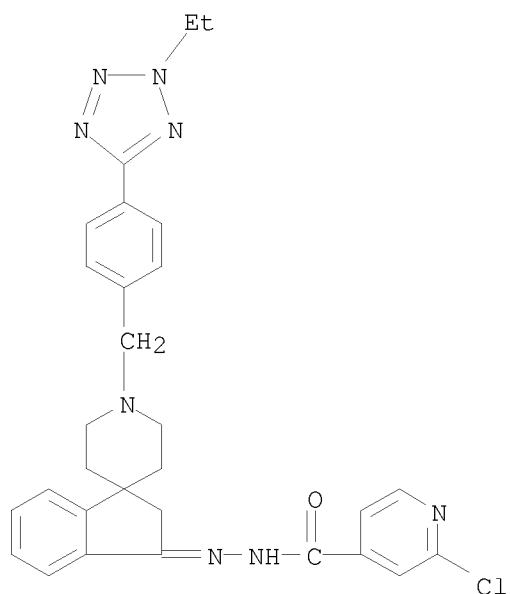


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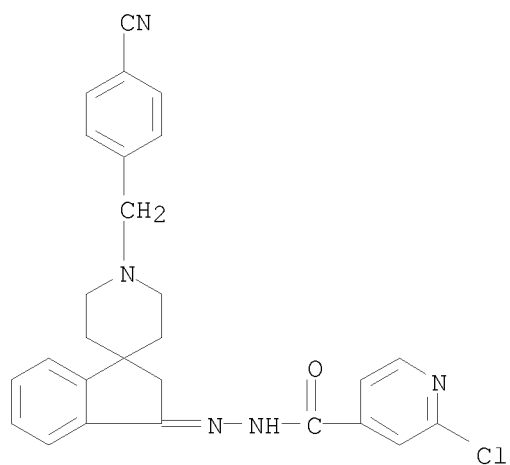
CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[1'-[[4-(2-ethyl-2H-tetrazol-5-yl)phenyl)methyl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

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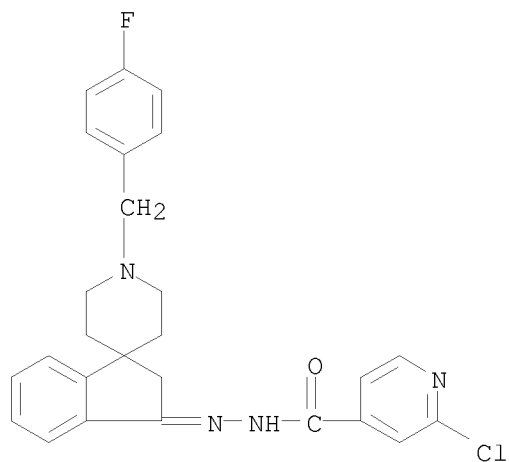
RN 855849-51-3 CAPLUS
CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[1'-[(4-cyanophenyl)methyl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-
ylidene]hydrazide (CA INDEX NAME)



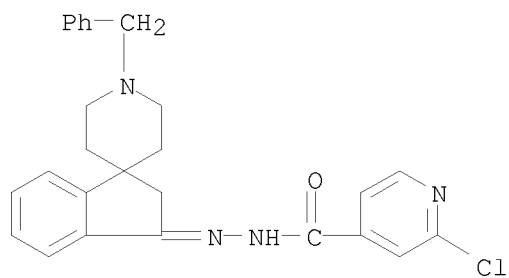
RN 855849-52-4 CAPLUS
CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[1'-[(4-fluorophenyl)methyl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-
ylidene]hydrazide (CA INDEX NAME)

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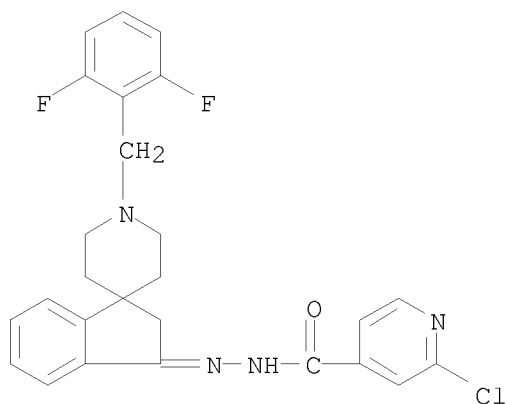


RN 855849-53-5 CAPLUS
CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[2,3-dihydro-1'-(phenylmethyl)spiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)



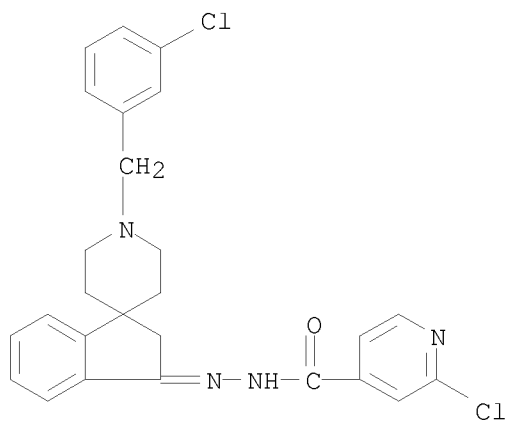
RN 855849-54-6 CAPLUS
CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[1'-[(2,6-difluorophenyl)methyl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

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RN 855849-55-7 CAPLUS

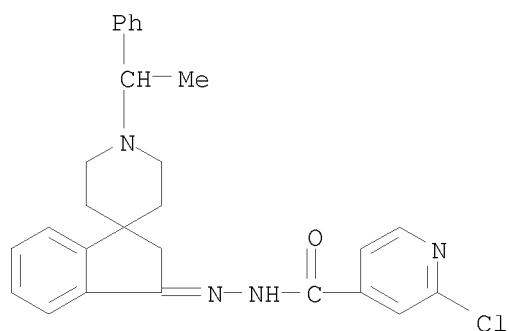
CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[1'-[(3-chlorophenyl)methyl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-
3-ylidene]hydrazide (CA INDEX NAME)



RN 855849-56-8 CAPLUS

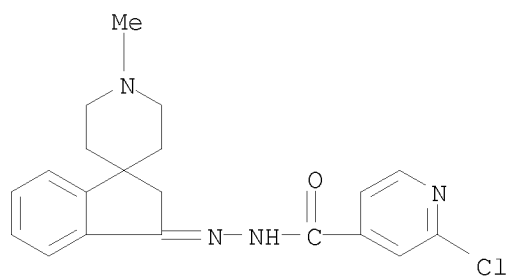
CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[2,3-dihydro-1'-(1-phenylethyl)spiro[1H-indene-1,4'-piperidin]-3-
ylidene]hydrazide (CA INDEX NAME)

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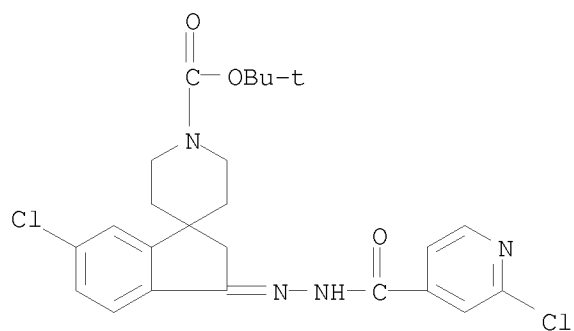
RN 855849-57-9 CAPLUS

CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-(2,3-dihydro-1'-methylspiro[1H-indene-1,4'-piperidin]-3-ylidene)hydrazide (CA INDEX NAME)



RN 855849-58-0 CAPLUS

CN Spiro[1H-indene-1,4'-piperidine]-1'-carboxylic acid,
6-chloro-3-[2-[(2-chloro-4-pyridinyl)carbonyl]hydrazinylidene]-2,3-dihydro-,
1,1-dimethylethyl ester (CA INDEX NAME)

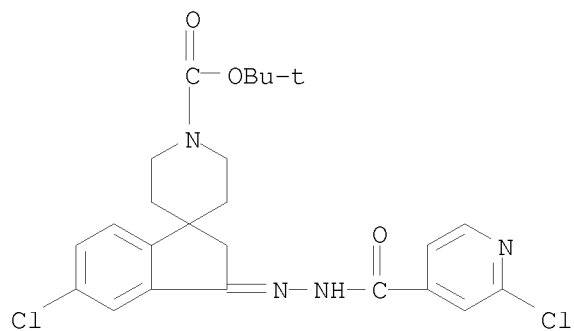


RN 855849-59-1 CAPLUS

CN Spiro[1H-indene-1,4'-piperidine]-1'-carboxylic acid,
5-chloro-3-[2-[(2-chloro-4-pyridinyl)carbonyl]hydrazinylidene]-2,3-dihydro-,
1,1-dimethylethyl ester (CA INDEX NAME)

26/08/2011

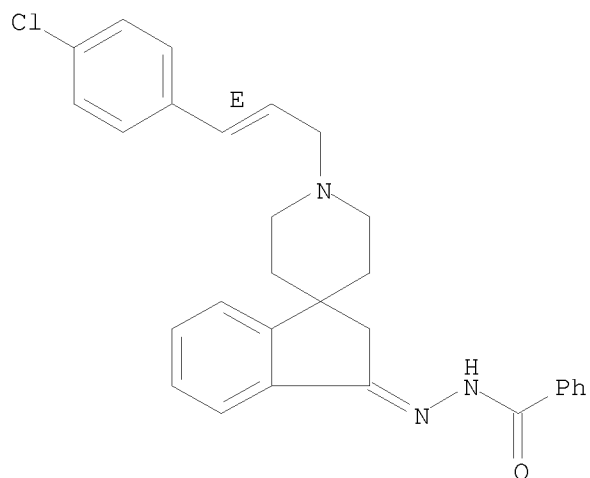
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RN 855849-60-4 CAPLUS

CN Benzoic acid, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

Double bond geometry as described by E or Z.

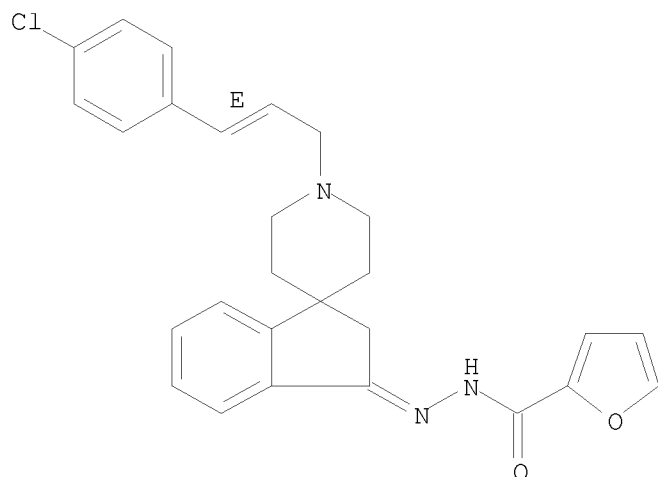


RN 855849-61-5 CAPLUS

CN 2-Furancarboxylic acid, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

Double bond geometry as described by E or Z.

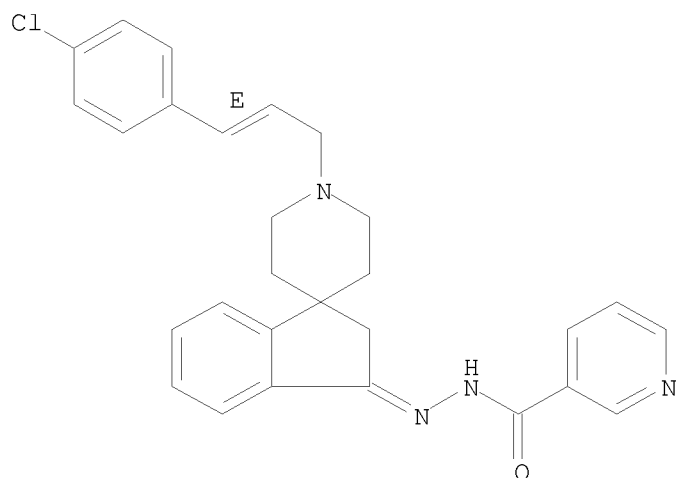
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RN 855849-62-6 CAPLUS

CN 3-Pyridinecarboxylic acid, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

Double bond geometry as described by E or Z.



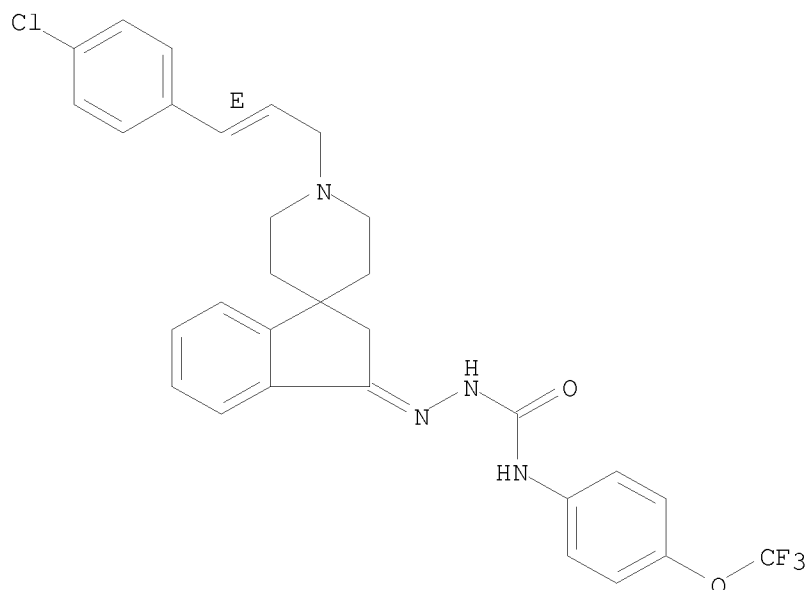
RN 855849-63-7 CAPLUS

CN Hydrazinecarboxamide, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]-N-[4-(trifluoromethoxy)phenyl] (CA INDEX NAME)

Double bond geometry as described by E or Z.

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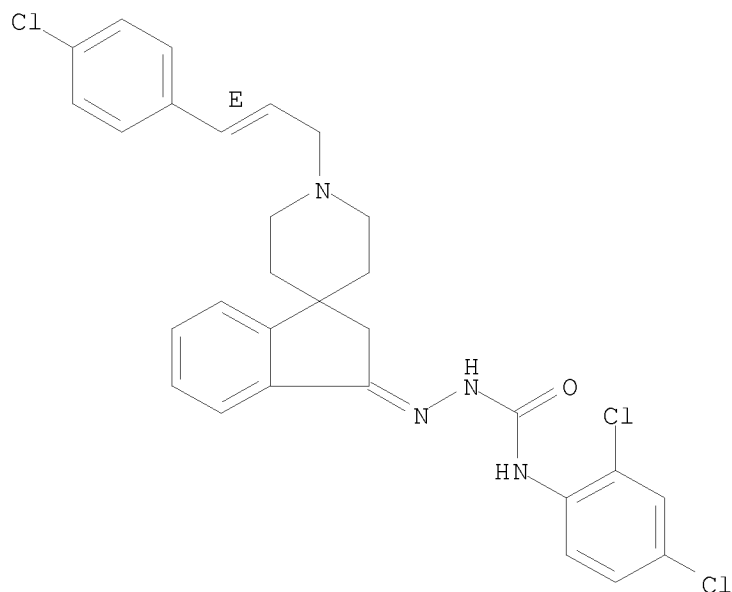
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RN 855849-64-8 CAPLUS

CN Hydrazinecarboxamide, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]-N-(2,4-dichlorophenyl)-(CA INDEX NAME)

Double bond geometry as described by E or Z.



RN 855849-65-9 CAPLUS

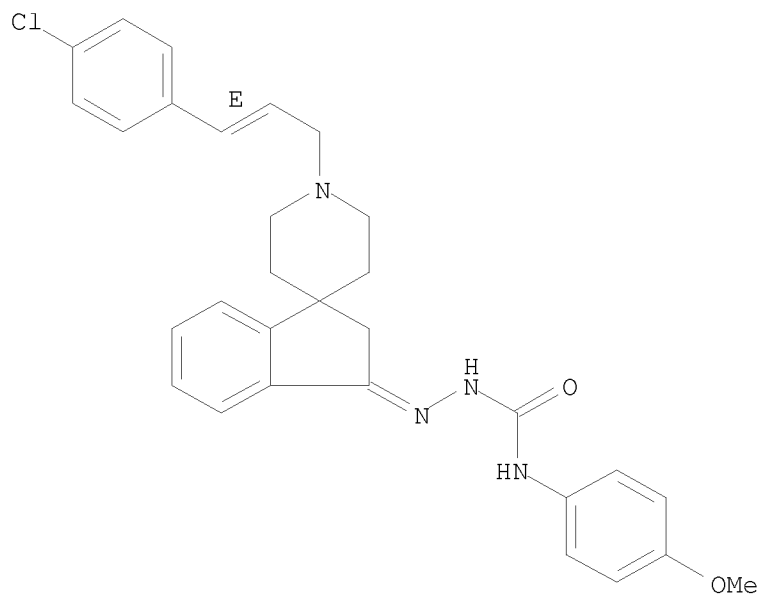
CN Hydrazinecarboxamide, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]-N-(4-methoxyphenyl)-

26/08/2011

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(CA INDEX NAME)

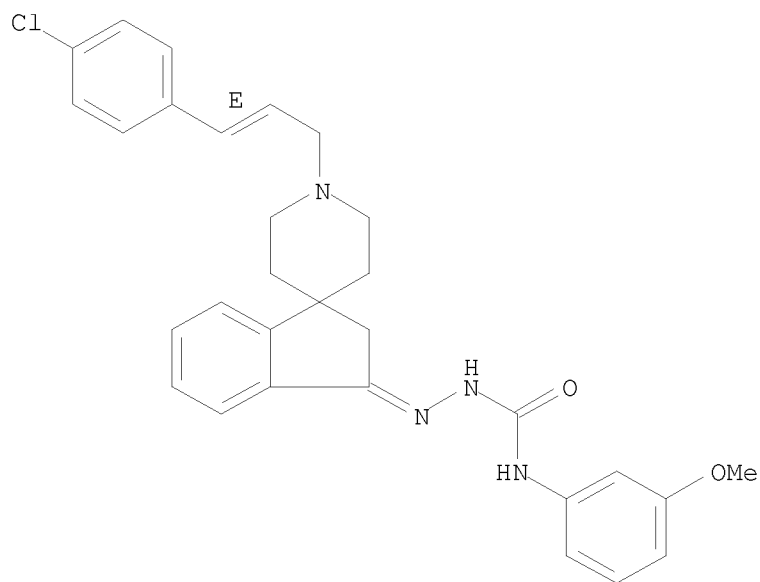
Double bond geometry as described by E or Z.



RN 855849-66-0 CAPLUS

CN Hydrazinecarboxamide, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]-N-(3-methoxyphenyl)-
(CA INDEX NAME)

Double bond geometry as described by E or Z.



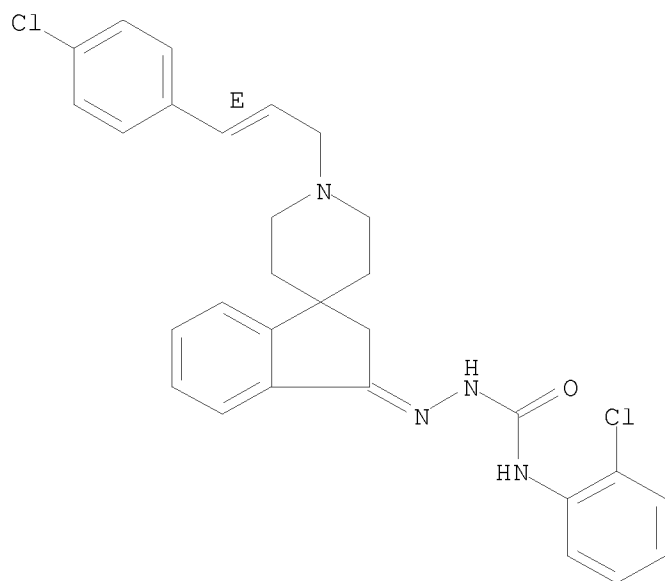
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26/08/2011

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CN Hydrazinecarboxamide, N-(2-chlorophenyl)-2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]- (CA INDEX NAME)

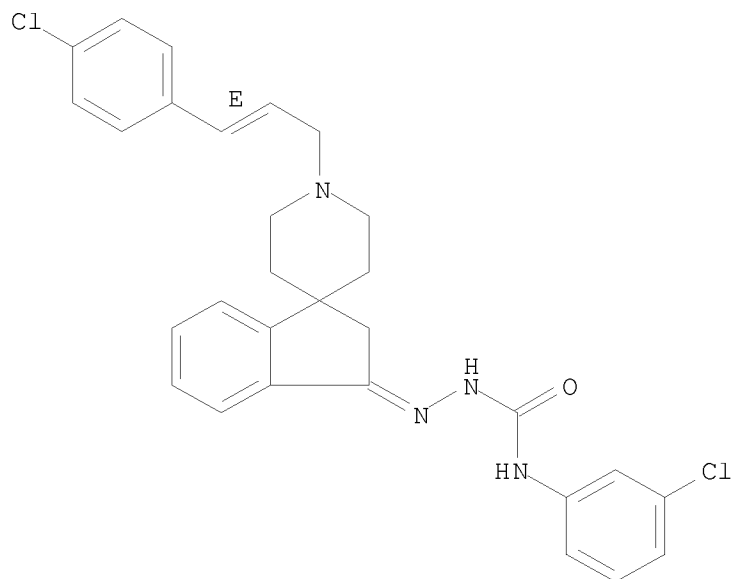
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RN 855849-68-2 CAPLUS

CN Hydrazinecarboxamide, N-(3-chlorophenyl)-2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]- (CA INDEX NAME)

Double bond geometry as described by E or Z.



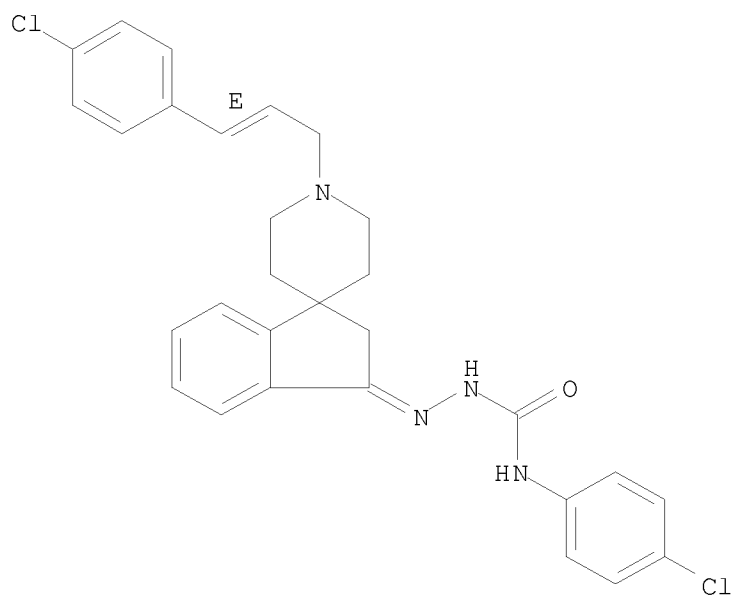
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RN 855849-69-3 CAPLUS

CN Hydrazinecarboxamide, N-(4-chlorophenyl)-2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]- (CA INDEX NAME)

Double bond geometry as described by E or Z.

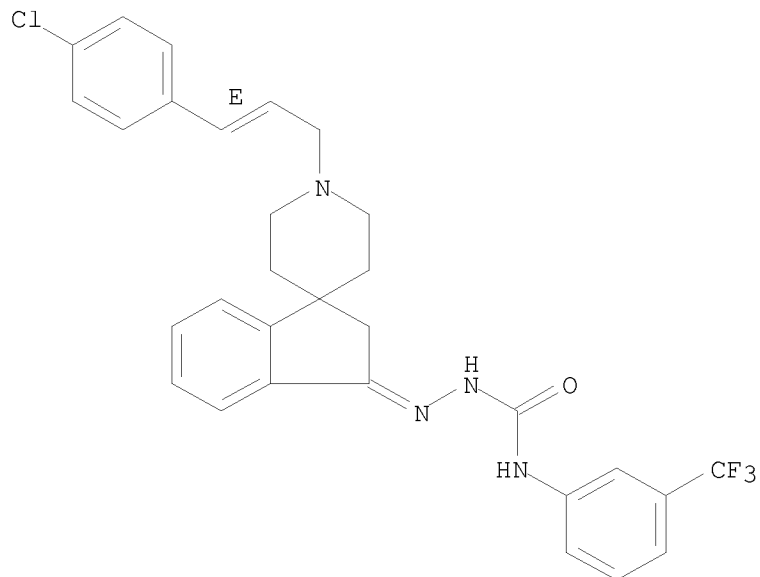


RN 855849-70-6 CAPLUS

CN Hydrazinecarboxamide, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)

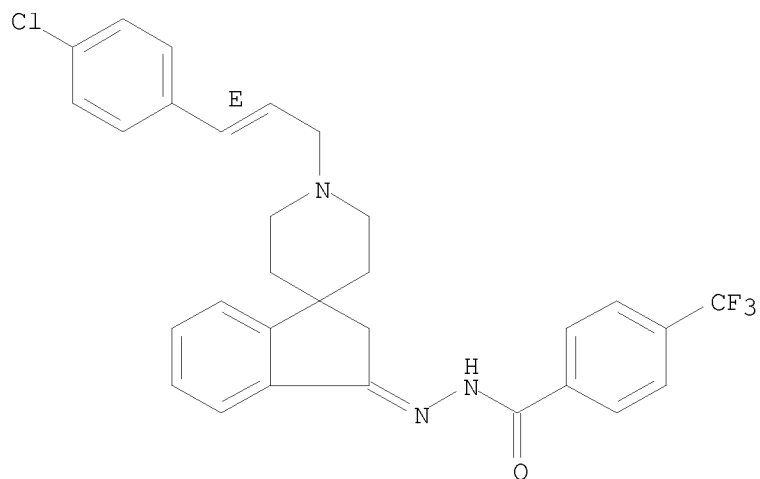
Double bond geometry as described by E or Z.

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RN 855849-71-7 CAPLUS
CN Benzoic acid, 4-(trifluoromethyl)-,
2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-
1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

Double bond geometry as described by E or Z.

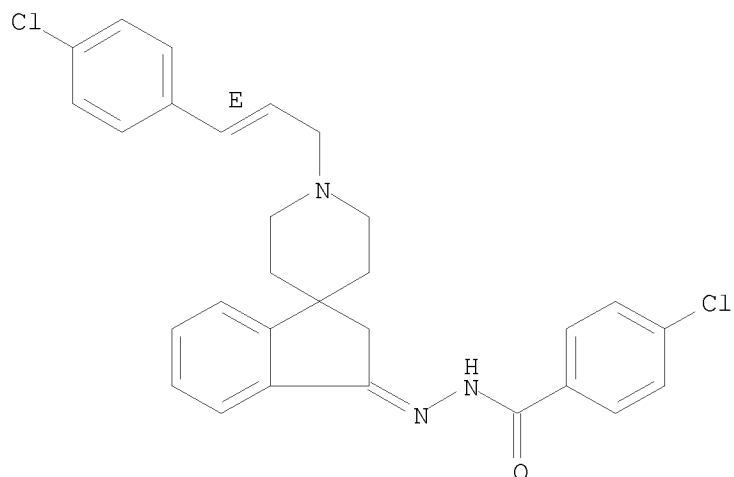


RN 855849-72-8 CAPLUS
CN Benzoic acid, 4-chloro-, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-
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NAME)

Double bond geometry as described by E or Z.

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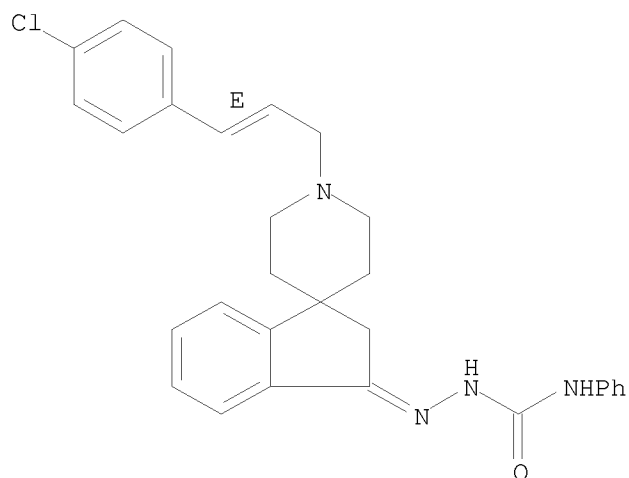
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RN 855849-73-9 CAPLUS

CN Hydrazinecarboxamide, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]-N-phenyl- (CA INDEX NAME)

Double bond geometry as described by E or Z.

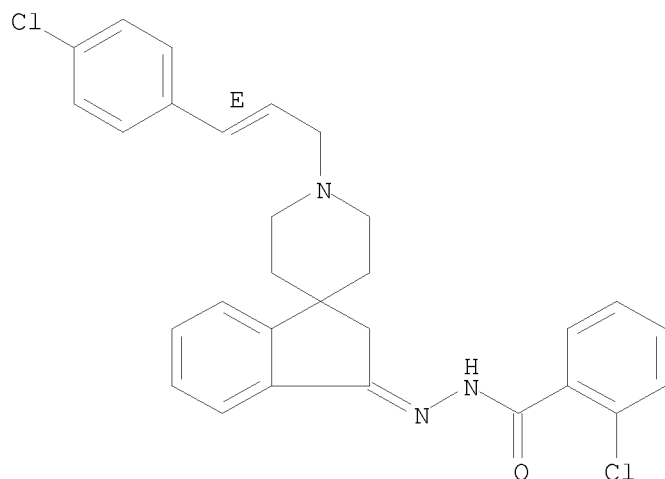


RN 855849-74-0 CAPLUS

CN Benzoic acid, 2-chloro-, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

Double bond geometry as described by E or Z.

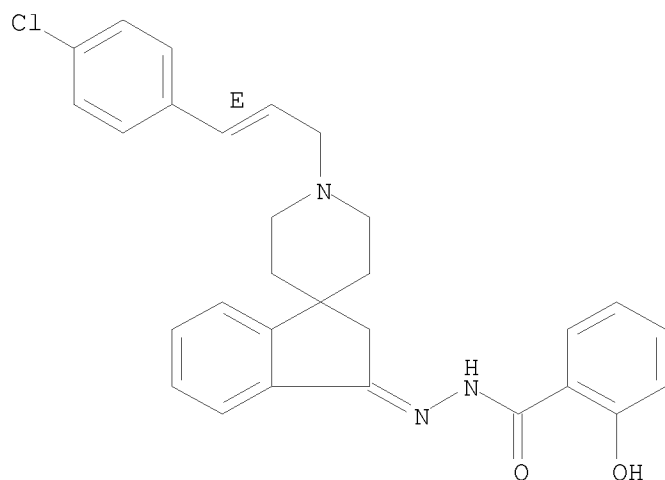
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CN Benzoic acid, 2-hydroxy-, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

Double bond geometry as described by E or Z.



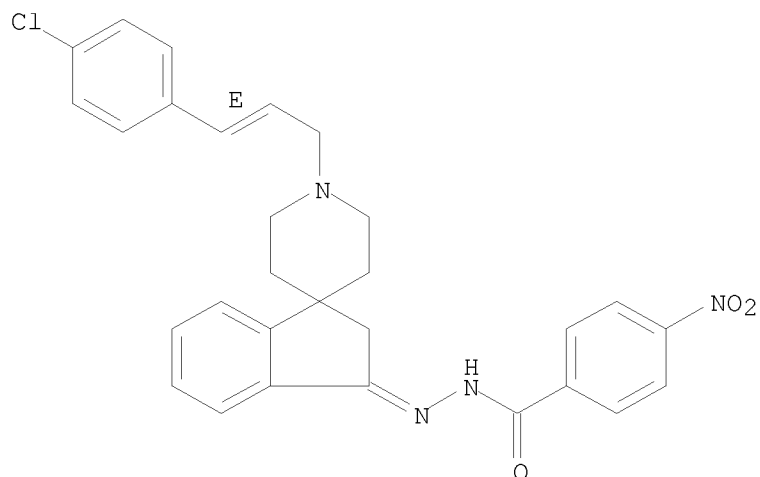
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CN Benzoic acid, 4-nitro-, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

Double bond geometry as described by E or Z.

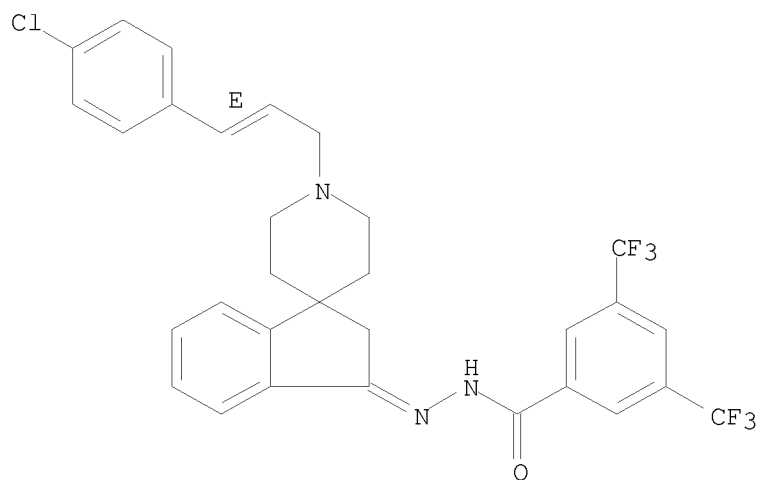
26/08/2011

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RN 855849-77-3 CAPLUS
CN Benzoic acid, 3,5-bis(trifluoromethyl)-, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

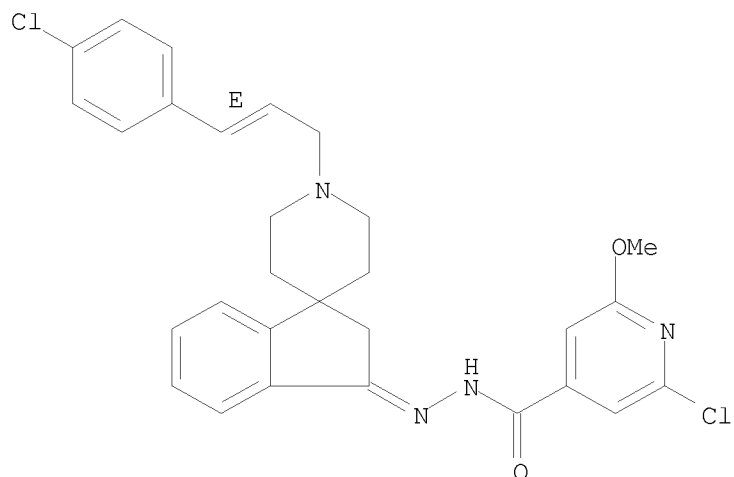
Double bond geometry as described by E or Z.



RN 855849-78-4 CAPLUS
CN 4-Pyridinecarboxylic acid, 2-chloro-6-methoxy-, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

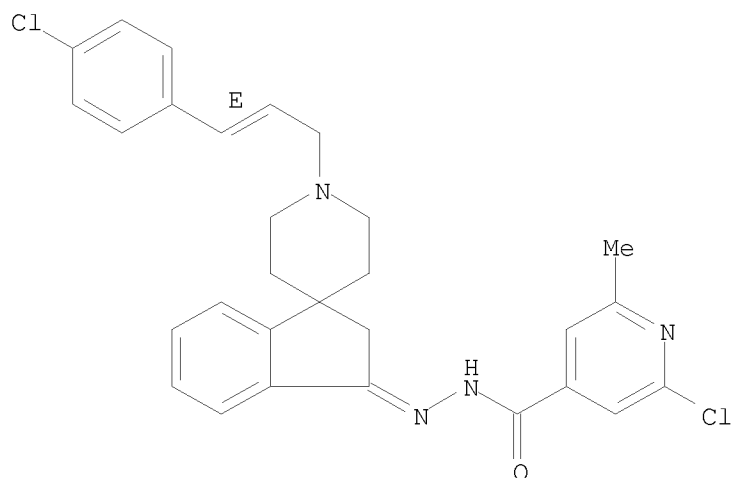
Double bond geometry as described by E or Z.

10581177.trn



RN 855849-79-5 CAPLUS
CN 4-Pyridinecarboxylic acid, 2-chloro-6-methyl-,
2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-
1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

Double bond geometry as described by E or Z.

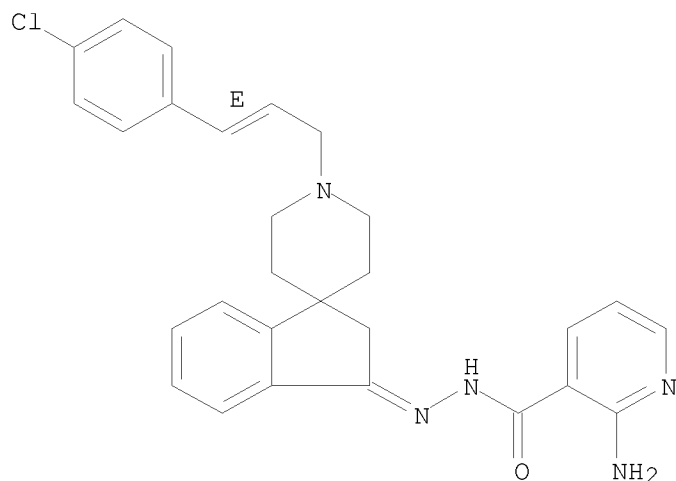


RN 855849-80-8 CAPLUS
CN 3-Pyridinecarboxylic acid, 2-amino-,
2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-
1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

Double bond geometry as described by E or Z.

26/08/2011

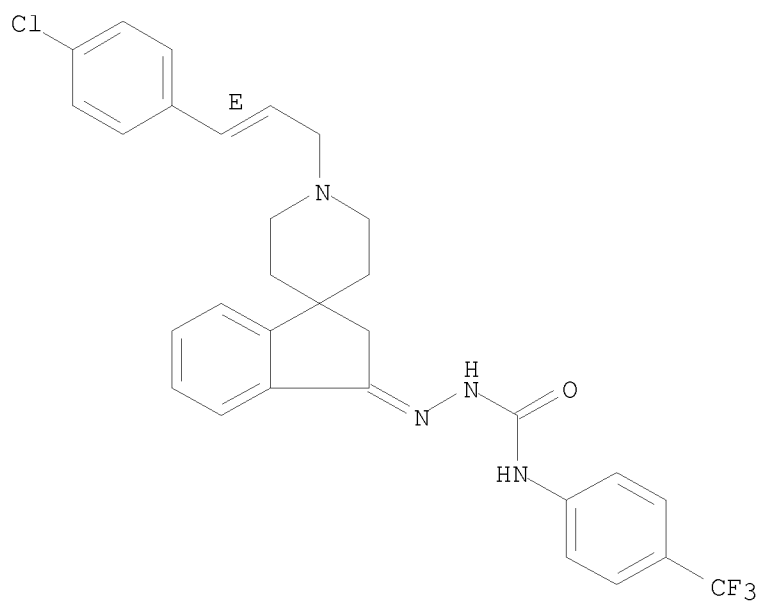
10581177.trn



RN 855849-81-9 CAPLUS

CN Hydrazinecarboxamide, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]-N-[4-(trifluoromethyl)phenyl]- (CA INDEX NAME)

Double bond geometry as described by E or Z.



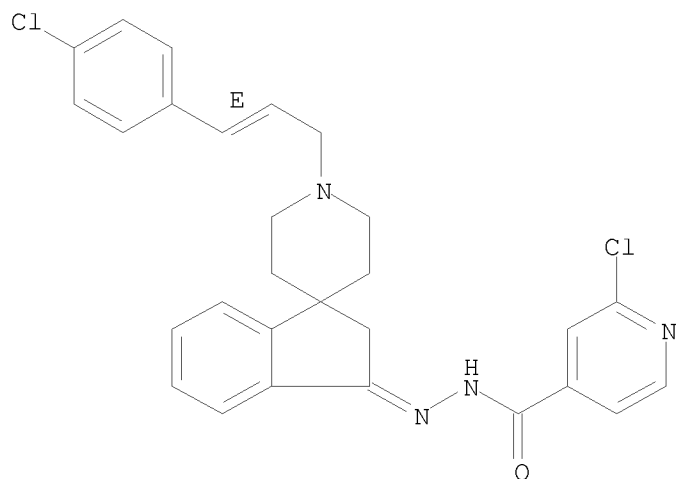
RN 855849-85-3 CAPLUS

CN 4-Pyridinecarboxylic acid, 2-chloro-, 2-[1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

Double bond geometry as described by E or Z.

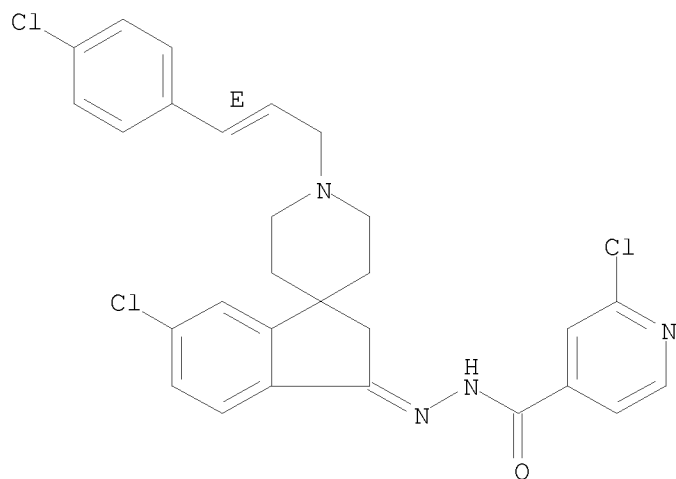
26/08/2011

10581177.trn



RN 855849-86-4 CAPLUS
CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[6-chloro-1'-[(2E)-3-(4-chlorophenyl)-2-propen-1-yl]-2,3-dihydrospiro[1H-
indene-1,4'-piperidin]-3-ylidene]hydrazide (CA INDEX NAME)

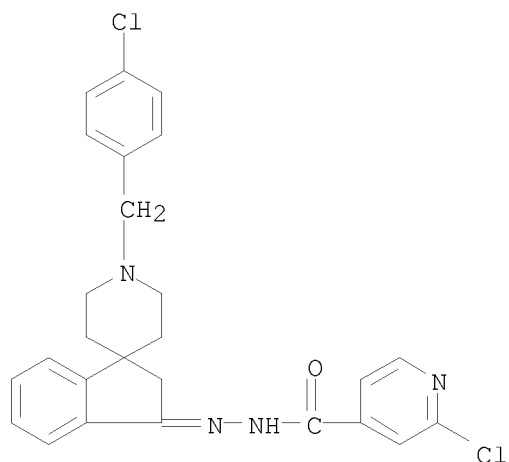
Double bond geometry as described by E or Z.



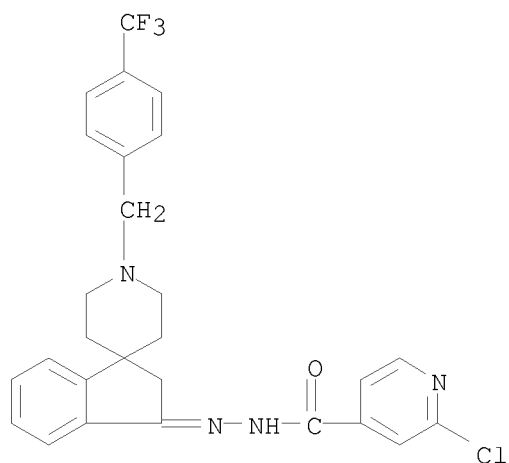
RN 855849-87-5 CAPLUS
CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[1'-[(4-chlorophenyl)methyl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-
3-ylidene]hydrazide (CA INDEX NAME)

26/08/2011

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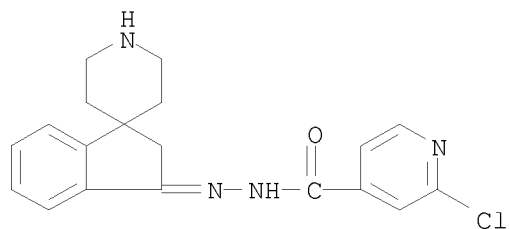
RN 855849-94-4 CAPLUS
CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-[2,3-dihydro-1'-[[4-(trifluoromethyl)phenyl]methyl]spiro[1H-indene-1,4'-
piperidin]-3-ylidene]hydrazide (CA INDEX NAME)



IT 855849-93-3P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT
(Reactant or reagent)
(preparation of insecticidal spiroindane derivs. as insecticides,
acaracides, molluscicides and nematocides)
RN 855849-93-3 CAPLUS
CN 4-Pyridinecarboxylic acid, 2-chloro-,
2-(2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-ylidene)hydrazide (CA
INDEX NAME)

26/08/2011

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RE.CNT 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

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=> file registry

C

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STRUCTURE FILE UPDATES: 25 AUG 2011 HIGHEST RN 1323485-64-8

DICTIONARY FILE UPDATES: 25 AUG 2011 HIGHEST RN 1323485-64-8

CAS Information Use Policies apply and are available at:

Uploading C:\Users\rdesai\Documents\STN Express 8.4\Queries\10581177.str



chain nodes :

15 16 17 18

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14

chain bonds :

8-15 16-17 17-18

ring bonds :

1-2 1-6 1-7 1-10 2-3 3-4 4-5 5-6 7-8 8-9 9-10 9-11 10-14 11-12 12-13
13-14

exact/norm bonds :

1-2 1-6 1-7 1-10 2-3 3-4 4-5 5-6 7-8 8-9 8-15 16-17 17-18

normalized bonds :

9-10 9-11 10-14 11-12 12-13 13-14

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom

11:Atom 12:Atom 13:Atom 14:Atom 15:CLASS 16:CLASS 17:CLASS 18:CLASS

L5 STRUCTURE UPLOADED

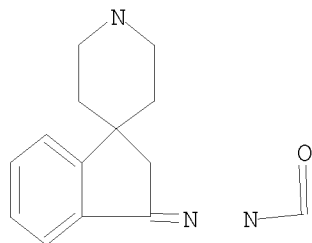
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L5 HAS NO ANSWERS

L5 STR

26/08/2011

10581177.trn



Structure attributes must be viewed using STN Express query preparation.

=> s 15

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SAMPLE SCREEN SEARCH COMPLETED - 1291 TO ITERATE

100.0% PROCESSED 1291 ITERATIONS

5 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 23665 TO 27975

PROJECTED ANSWERS: 5 TO 234

L6 5 SEA SSS SAM L5

=> s 15 ful

FULL SEARCH INITIATED 16:20:06 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 26312 TO ITERATE

100.0% PROCESSED 26312 ITERATIONS

42 ANSWERS

SEARCH TIME: 00.00.01

L7 42 SEA SSS FUL L5

C

CAlplus now includes complete International Patent Classification (IPC) reclassification data for the second quarter of 2011.

CAS Information Use Policies apply and are available at:

<http://www.cas.org/legal/infopolicy.html>

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 17

L8 3 L7

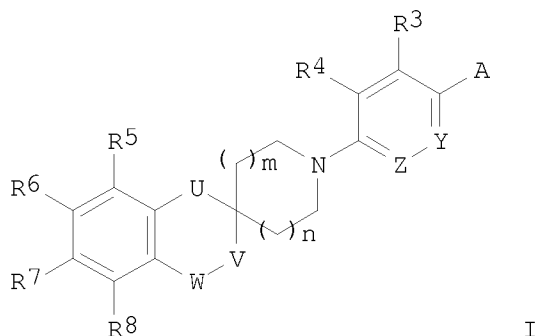
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26/08/2011

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26/08/2011

L8 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2011 ACS on STN
GI



AB Disclosed is an agent represented by a general formula I (A = CONHR₁, ER₂; R₁ = (un)substituted C1-6 alkyl, (un)substituted C3-6 cycloalkyl; E = 1,3,4-oxadiazol, 1,2,4-oxadiazol, imidazol; R₂ = H, (un)substituted C1-6 alkyl, phenyloxy, phenylamino, pyridylamino; R₃, R₄ = H, halogen, C1-6 alkyl, R₅, R₆, R₇, R₈ = H, halogen, C1-6 (halogenated)alkyl, C1-6 (halogenated)alkoxy; Z, Y = N, :CH- except Z = Y = :CH-; U = single bond, methylene, CH(OH), carbonyl, O, S, sulfinyl, sulfonyl; V = single bond, methylene, CH(OH); W = methylene, CH(OH), etc.; m, n = 0, 1 except m = 0/n = 1 and m = 1/n = 0), or is pharmaceutically acceptable salt as an active component. The agent has stearyl-CoA desaturase-inhibitory effect, and suitable for use for prevention and/or treatment of related disease, e.g. obesity, hyperlipidemia, lipid metabolism disorder, diabetes, etc. For example, 6-(3,4-dihydro-1'H-spiro[chromene-2,4'-piperidin]-1'-yl)-N-(2-hydroxy-2-phenylethyl)piridazine-3-carboxamide (II) was prepared. The compound II showed ≥ 50 % stearyl-CoA desaturase activity-inhibitory effect at 0.2 μM in vitro. Also, a capsule containing II 50 mg/150 mg capsule was formulated.

AN 2009:1430440 CAPLUS

DN 151:565181

TI Pharmaceutical agent containing novel spiro piperidine derivatives having stearyl-CoA desaturase-inhibitory effect

IN Uto, Yoshikazu; Kiyotsuka, Yohei

PA Daiichi Sankyo Co., Ltd., Japan

SO Jpn. Kokai Tokkyo Koho, 136 pp.

CODEN: JKXXAF

DT Patent

LA Japanese

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 2009269850	A	20091119	JP 2008-120937	20080507
PRAI	JP 2008-120937		20080507		
OS	MARPAT 151:565181				
IT	1024604-69-0P				

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

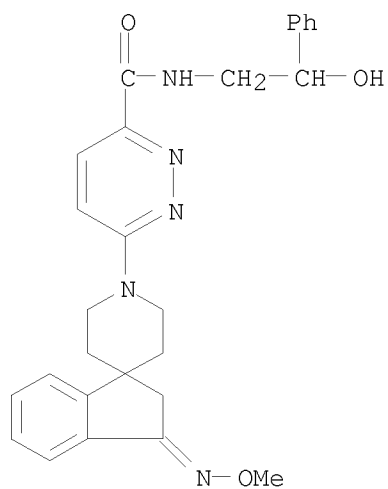
(novel spiro piperidine derivs. having stearyl-CoA

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desaturase-inhibitory effect)

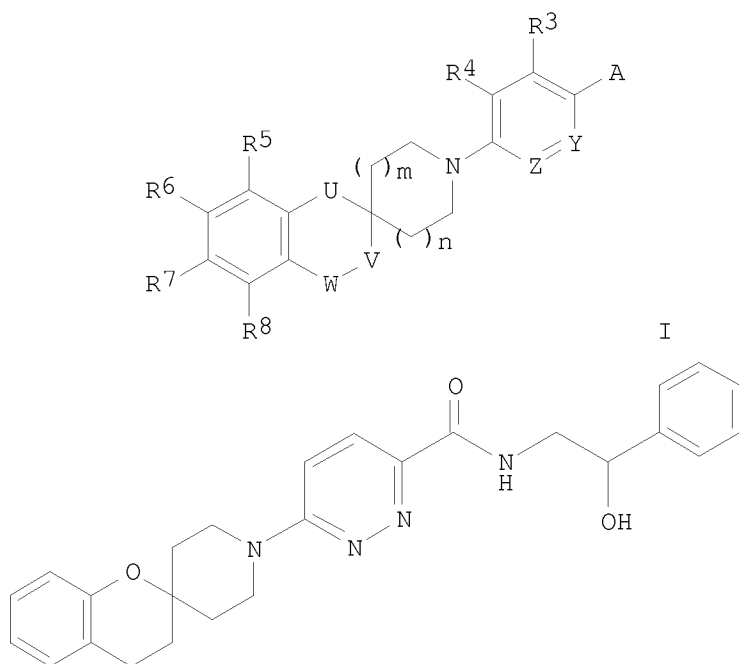
RN 1024604-69-0 CAPLUS

CN 3-Pyridazinecarboxamide, 6-[2,3-dihydro-3-(methoxyimino)spiro[1H-indene-1,4'-piperidin]-1'-yl]-N-(2-hydroxy-2-phenylethyl)- (CA INDEX NAME)



OSC.G 1 THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD (1 CITINGS)

L8 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2011 ACS on STN
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II

AB The title compds. [I; A = CONHR₁ or ER₂; R₁ = (un)substituted C1-6 alkyl or C3-6 cycloalkyl; E = 1,3,4-oxadiazole, 1,2,4-oxadiazole, or imidazole group; R₂ = H, each (un)substituted C1-6 alkyl, phenyloxy, phenylamino, or pyridylamino; R₃, R₄ = H, halo, C1-6 alkyl; R₅-R₈ = H, halo, C1-6 alkyl, C1-6 haloalkyl, C1-6 alkoxy, C1-6 haloalkoxy; Z, Y = N, CH; U = single bond, methylene, CH(OH), CH(OQ₁), CO, C(:NOQ₂), O; Q₁, Q₂ = C1-6 alkyl; m, n = 0 or 1, provided that a case where m = 0 and n = 1 or m = 1 and n = 0 is excluded] or pharmacol. acceptable salts were prepared. These compds. have excellent stearoyl-CoA-desaturase inhibitory activity and are useful for the prevention and/or treatment of obesity, hyperlipidemia, hypertriglyceridemia, insulin resistance syndrome, abnormal glucose tolerance, diabetes, diabetes complications (including diabetes peripheral nerve disorder, diabetic nephropathy, diabetic retinopathy, and diabetic macroangiopathy), cataract, gestational diabetes, polycystic ovary syndrome, arteriosclerosis, atherosclerosis, diabetic arteriosclerosis, hypertension, cerebral vascular disorders, coronary artery disease, fatty liver, nonalcoholic fatty hepatitis, dyspnea, backache (lumbago), gonarthrosis, gout, and cholelithiasis. Thus, a suspension of 92 mg 6-chloro-N-(2-hydroxy-2-phenylethyl)pyridazine-3-carboxamide, 67 mg 3,4-Dihydro-1'H-spiro[chromene-2,4'-piperidine], and diisopropylethylamine in n-butanol was heated at 120° for 25 h to give 49% 6-(3,4-dihydrospiro[chromene-2,4'-piperidin]-1'-yl)-N-(2-hydroxy-2-phenylethyl)pyridazine-3-carboxamide (II). II at 2 μM inhibited ≥50% stearoyl-CoA-desaturase. A capsule and a tablet formulation containing II were described.

10581177.trn

AN 2008:583341 CAPLUS
DN 148:538285
TI Preparation of spiropiperidine derivatives as stearyl-CoA-desaturase
inhibitors
IN Uto, Yoshikazu; Kiyotsuka, Yohei
PA Daiichi Sankyo Company, Limited, Japan
SO PCT Int. Appl., 233pp.
CODEN: PIXXD2
DT Patent
LA Japanese
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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PI	WO 2008056687	A1	20080515	WO 2007-JP71605	20071107
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BH, BR, BW, BY, BZ, CA,				
	CH, CN, CO, CR, CU, CZ, DE, DK, DM, DO, DZ, EC, EE, EG, ES, FI,				
	GB, GD, GE, GH, GM, GT, HN, HR, HU, ID, IL, IN, IS, JP, KE, KG,				
	KM, KN, KP, KR, KZ, LA, LC, LK, LR, LS, LT, LU, LY, MA, MD, ME,				
	MG, MK, MN, MW, MX, MY, MZ, NA, NG, NI, NO, NZ, OM, PG, PH, PL,				
	PT, RO, RS, RU, SC, SD, SE, SG, SK, SL, SM, SV, SY, TJ, TM, TN,				
	TR, TT, TZ, UA, UG, US, UZ, VC, VN, ZA, ZM, ZW				
	RW: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE,				
	IS, IT, LT, LU, LV, MC, MT, NL, PL, PT, RO, SE, SI, SK, TR, BF,				
	BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG, BW,				
	GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ,				
	BY, KG, KZ, MD, RU, TJ, TM				

PRAI JP 2006-303866 A 20061109

OS MARPAT 148:538285

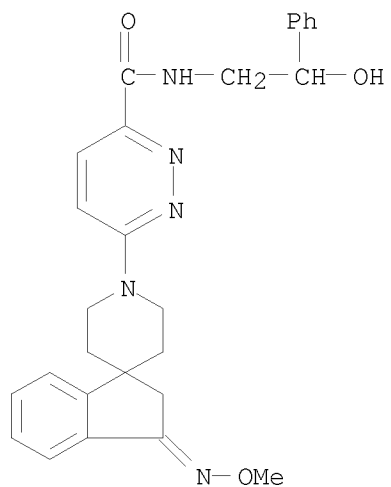
IT 1024604-69-0P, N-(2-Hydroxy-2-phenylethyl)-6-[3-(methoxyimino)-
2,3-dihydrospiro[indene-1,4'-piperidin]-1'-yl]pyridazine-3-carboxamide
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
(Therapeutic use); BIOL (Biological study); PREP (Preparation); USES
(Uses)

(preparation of spiropiperidine derivs. as stearyl-CoA-desaturase
inhibitors)

RN 1024604-69-0 CAPLUS

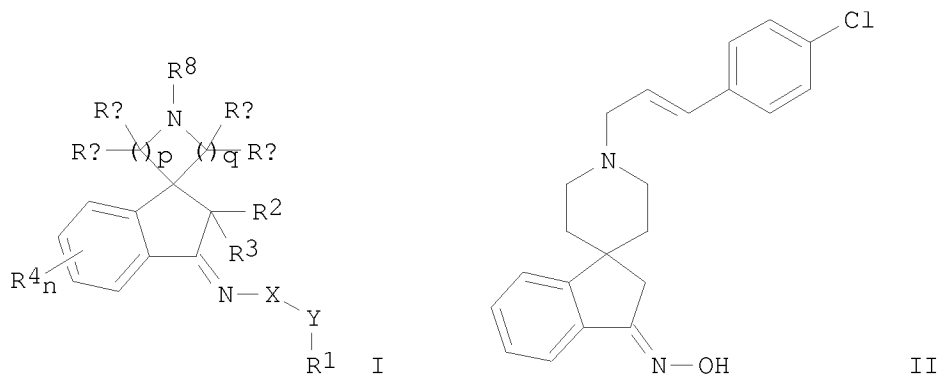
CN 3-Pyridazinecarboxamide, 6-[2,3-dihydro-3-(methoxyimino)spiro[1H-indene-
1,4'-piperidin]-1'-yl]-N-(2-hydroxy-2-phenylethyl)- (CA INDEX NAME)

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OSC.G	7	THERE ARE 7 CAPLUS RECORDS THAT CITE THIS RECORD (7 CITINGS)
RE.CNT	34	THERE ARE 34 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT		

L8 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2011 ACS on STN
GI



AB Title compds. I [X = O, amino; Y = bond, CO, CS, SO0-2; R1 = H, alkyl, alkoxy carbonyl, etc.; R2-3 = H, halo, CN, alkyl, etc.; R4 = halo, NO2, CN, etc.; Ra = H, halo, OH, CN, etc.; p, q = 0-6; R8 = alk(en/yn)yl, etc.] are prepared For instance, II is prepared in 3 steps from spiro[indan-1-one-3,4'-piperidine]-1'-carboxylic acid tert Bu ester, 4-chlorocinnamyl chloride and hydroxylamine (E (dominant) and Z oximes isolated). Selected example compds. gave >80% control of *Spodoptera littoralis*. I are useful in controlling insects, acarines, nematodes or molluscs.

AN 2005:570877 CAPLUS

DN 143:77964

TI Preparation of insecticidal spiroindane derivatives

IN Cassayre, Jerome; Molleyres, Louis-Pierre; Maienfisch, Peter; Cederbaum, Fredrik

PA Syngenta Participations A.-G., Switz.

SO PCT Int. Appl., 114 pp.

CODEN: PIXXD2

DT Patent

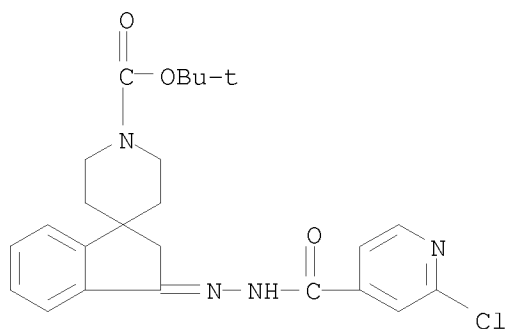
LA English

FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005058836	A1	20050630	WO 2004-IB4108	20041209
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RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
EP 1697327	A1	20060906	EP 2004-806338	20041209
EP 1697327	B1	20110713		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,				

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IE, SI, LT, FI, RO, CY, TR, BG, CZ, EE, HU, PL, SK, IS
BR 2004017555 A 20070327 BR 2004-17555 20041209
JP 2007516253 T 20070621 JP 2006-543659 20041209
AT 516273 T 20110715 AT 2004-806338 20041209
IN 2006CN02077 A 20070706 IN 2006-CN2077 20060612
US 20080306101 A1 20081211 US 2008-581177 20080828
PRAI GB 2003-28906 A 20031212
WO 2004-IB4108 W 20041209
ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT
OS CASREACT 143:77964; MARPAT 143:77964
IT 855849-47-7P
RL: AGR (Agricultural use); BSU (Biological study, unclassified); RCT
(Reactant); SPN (Synthetic preparation); BIOL (Biological study); PREP
(Preparation); RACT (Reactant or reagent); USES (Uses)
(preparation of insecticidal spiroindane derivs. as insecticides,
acaracides, molluscicides and nematocides)
RN 855849-47-7 CAPLUS
CN Spiro[1H-indene-1,4'-piperidine]-1'-carboxylic acid,
3-[2-[(2-chloro-4-pyridinyl)carbonyl]hydrazinylidene]-2,3-dihydro-,
1,1-dimethylethyl ester (CA INDEX NAME)



RE.CNT 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

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COST IN U.S. DOLLARS

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